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244389

SUPERFUND TECHNICAL ASSESSMENT AND RESPONSE TEAM EPA CONTRACT 68-W5-0019

START-02-F-01540

TRANSMITTAL MEMO

To:

Jeff Bechtel, OSC

Removal Action Branch, U.S. EPA Region II

From:

Smita Sumbaly, Data Reviewer

START Region II

Subject:

Color Technology Site

Data Validation Assessment

Date:

December 17, 1997.

The purpose of this memo is to transmit the following information:

• Data validation results for the following parameters:

TAL metals

11 samples

Full TCL

21 samples

Matrices and Number of Samples

Soil

11 samples

Wipe

10 samples

Sampling date:

September 19, 1997.

The final data assessment narrative and original analytical data package are attached.

cc:

START PM

Michael Mahnkopf

START FILE

TDD #:02-97-09-0013

TDD #:02-97-05-0021

PCS #:2117

U.S. ENVIRONMENTAL PROTECTION AGENCY

MEMORANDUN	1	, <u> </u>					
DATE:	December 16	i, 1997	•		.*		
то:	Jeff Bechtel, USEPA Regi						
FROM:	Smita Sumb		m				
SUBJECT:	QA/QC Con	npliance Rev	iew Summa	ry			• .
Sp Su Ma	npared to EPA st	andards for co	Blanks DFTPP a Chromat Holding	Measures for and BFB Tu ography	or the following	ted have b owing gen	een eral
	asures used to sup	port the follow	wing conclus	sions are att	ached so	that the rev	/iew
<u>Summary</u>	of Results	I	II	III Post	[/ PCBs	IV	
Acceptable as Su Acceptable with Unacceptable, A Unacceptable	Comments	X	X	X	——————————————————————————————————————		
Data Reviewed b	oy:	Am.	495	mbeh	Date:_	12/10	<u>5/9</u> 7
Approved By:		m Soi	r	. .	Date:_	14/1/4	7

Area Code/Phone No.:

NARRATIVE

CASE No. 2117

SITE NAME:	Color Technology Site	<u> </u>		-
	60 Cornell Blvd, Some	rville, New Jersey.		_
Laboratory Name:	Industrial Corrosion M	anagement, Inc.(ICM))	
INTRODUCTION:				
The laboratory's porti	on of this Case consisted	of 21 samples collec	ted on Septemb	er 19, 1997.
	ted problem(s) with the	And the second s		11
Lab did not perfo hexane preservative.	rmed volatile analysis o	n wipe samples due	to analytical pr	oblems with
The laboratory re TCL(VOA/BNA/Pes	ported <u>No</u> proble st/PCBs) - Organic Comp	ms with the analy oounds.	ses of sample	es for Full
have been accessed	mmented on the criteria but no discussion is given branch or require no commed.	en where the evaluator	r has determined	u mai cinena

Evaluation by Fraction

I. Volatile (VOAs)

Y Holding Time Y MS/MSD

Y GC/MS Tuning Y Compound ID (HSL, TIC)

Y Calibration, Initial Y Spectra Quality

Y Calibration, Continuing Y Standards

Y Blank Y Chromatography

Y Surrogate Recovery Y Data Completeness

Y Laboratory Fortified Blank Y Laboratory Storage Blank

Comments:

1. Refer to Data Assessment Narrative.

II. Base Neutral/Acids (BNAs)

Y Holding Time Y MS/MSD

Y GC/MS Tuning Y Compound ID (HSL, TIC)

Y Calibration, Initial Y Spectra Quality

Y Calibration, Continuing Y Standards

Y Blank Y Chromatography

Y Surrogate Recovery Y Data Completeness

Comments:

1. Refer to Data Assessment Narrative.

III. Pesticides

Y Holding Times Y Calibration Linearity

Y Instrument Performance Y Blank

Y Surrogate Recovery Y Retention Time Window

Y MS/MSD Y Analytical Sequence

Y Compound ID (HSL,TIC) Y RT Check for TCX

Y Chromatography

Comments:

1. Refer to Data Assessment Narrative.

REGION II START DATA ASSESSMENT REPORT

RFP Project #: 2117	Case #: NA		SDG #: <u>CTS01</u>	
LAB: Industrial Corrosion	Management, Inc.	LAB Code:	<u>ICM</u>	
SITE: Color Technology		Analysis: To	CL Organic Parameters Ma	trix:
Contractor: START	Reviewer: Smita Sun	nbaly	Water Soil/Sedime	
CERCLIS ID #:			Wij	pe:09_
The current Functional Gu	idelines for evaluating or	ganic data have	been applied.	
(presumptive evidence for the				4
Two facts should be noted unusable. In other words information as to whether the because they cannot be reliable.	, due to significant QC parties to significant QC parties to significant QC parties are significant quality and parties are significant quality	problems, the a not. "R" value resort. The seco	naiysis is invalid and j s should not appear on and fact to keep in min	data table d is that n
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On 19 September 1997, USEPA Region II - personnel collected 11 surface soil and 10 wipe samples for Target Compound List (TCL) organic analyses from the Color Technology Site, Somerville, New Jersey. Within twenty-four hours of collection, samples were picked up by Industrial Corrosion Management (ICM), Randolph, New Jersey. The laboratory verified that samples were received intact, properly custody sealed, and refrigerated (sample cooler temperature recorded at 2.8°C).

Target Compound List (TCL) organic analyses were performed following the Contract Laboratory Program (CLP) Statement of Work (SOW) number OLM03.1.

Client identification (ID) and laboratory ID numbers:

Client ID No.	Laboratory ID No.	<u>Matrix</u>
CTS01	272610	Soil
CTS02	272611	Soil
CTS03	272612	Soil
CTS04	272613	Soil
CTS05	272614	Soil
CTS06	272615	Soil
CTS07	272616	Soil
CTS08	272617	Soil
CTS09	272618	Soil
CTS010	272619	Soil
CTS011 1	272620	Soil
CTW01	272621	Wipe
CTW02	272622	Wipe
CTW03	272623	Wipe
CTW04	272624	Wipe
CTW05	272625	Wipe
CTW06	272626	Wipe
CTW07	272627	Wipe
CTW08	272628	Wipe
CTW09 ²	272629	Wipe
CTW10(BLANK)	272630	Wipe

Sample CTS011 is a field duplicate sample of sample CTS01.

Sample CTW09 is a field duplicate sample of sample CTW01.

1. HOLDING TIMES:

The amount of an analyte in a sample can change with time due to chemical instability, degradation, volatilization, etc. If the specified holding time is exceeded, the data may not be valid. Those analytes detected in the samples whose holding time has been exceeded will be qualified as estimated, "J". The non-detects (sample quantitation limits) will be flagged as estimated, "J", or unusable, "R", if the holding times are grossly exceeded.

The following analytes in the samples shown were qualified because of holding time:

TCL Data

<u>VOA</u> - The following data were qualified as estimated "J" or rejected "R" due to exceeding holding time criteria:

Sample ID	<u>Matrix</u>	Date Sampled	<u>Date</u> <u>Analyzed</u>	<u>Oualifier</u>	# Compounds
CTS05	Soil	09/19/97	09/29/97	"J"	33

Note: If properly preserved, aqueous samples maintained at 4°C must be analyzed within fourteen (14) days of collection. If unpreserved, aqueous samples must be analyzed within seven (7) days for aromatic hydrocarbons. Soil/Solid samples must be analyzed within ten (10) days of collection.

BNA - The following data were qualified as estimated "J" or rejected "R" due to exceeding holding time criteria:

Sample ID Matrix Date Sampled Date Extracted Date Analyzed Qualifier # Compounds

Data met the QC criteria

<u>Pest/PCBs</u> - The following data were qualified as estimated "J" or rejected "R" due to exceeding holding time criteria:

Sample ID Matrix Date Sampled Date Extracted Date Analyzed Qualifier # Compounds

Data met the QC criteria

Note: Continuous extraction of water samples must be started within seven (7) days of the date of collection. Soil/Sediment/Solid samples must be extracted within ten (10) days of collection. Extracts must be analyzed within forty (40) days of extraction.

2. BLANK CONTAMINATION:

Quality Assurance (QA) blanks [i.e., method, trip, field or rinse blanks] are prepared to identify any contamination which may have been introduced into the samples during sample preparation or field activity. Method blanks measure laboratory contamination. Trip blanks measure cross-contamination of samples during shipment. Field and rinse blanks measure cross-contamination of samples during field operations. If the concentration of the analyte is less than 5 times the blank contaminant level (10 times for common contaminants), the analytes are qualified as non-detects, "U". The following analytes in the samples shown were qualified with "U" for these reasons:

A) Method Blank Contamination

<u>VOA</u> - The following compounds were qualified as non-detected "U" in the associated samples due to method blank contamination:

Compound

Associated Samples

Methylene Chloride

CTS-01, -03, -06, -08, -09, -10Re, -11, -07Re & -02

<u>VOA</u> - The following TICs were rejected "R" in the indicated samples due to detection in the associated method blank:

<u>TIC</u>	Associated Method Blank	Associated Samples
Column Bleed	VBLKSB VBLKSC VBLKSD	CTS-01 & CTS-06 CTS-04, -06Re, -08Re, -09Re & 10 CTS-08, -09, -10Re, -11, -07Re, -02 & -07

BNA - The following compounds were qualified as non-detected "U" in the associated samples due to method blank contamination:

Compound

Associated Samples

bis(2-Ethylhexyl)phthalate

CTS-04, -04DL, -06, -06DL & -02

BNA - The following TICs were rejected "R" in the indicated samples due to detection in the associated method blank:

TIC

Associated Method Blank

Associated Samples

Ethanol, 2-(2-ethoxyethoxy)-

SBLK26

CTS-10, -05 & -07

<u>Pest/PCBs</u> - The following compounds were qualified as non-detected "U" in the associated samples due to method blank contamination:

Data met the QC criteria

B) Field or Rinse Blank Contamination ("water blanks" or "distilled water blanks" are validated like any other sample)

<u>Full TCL</u> - The following compounds were qualified as non-detected "U" in the associated samples due to rinse blank contamination:

Compound

Associated Samples

Not Applicable

C) Trip Blank Contamination

<u>VOA</u> - The following compounds were qualified as non-detected "U" in the associated samples due to trip blank contamination:

Compound

Associated Samples

Not Applicable

D) TIC Blank Contamination

<u>VOA/BNA</u> - The following TIC contaminants were rejected "R" in the indicated samples because these compounds are target TCL compounds also present in the associated volatile or semi-volatile fraction analyses:

Fraction	TIC	Samples
BNA	Fluorene Benzo (k) Fluoranthene Pyrene	CTS-04 CTS-04DL CTS-06

<u>VOA/BNA</u> - All TICs qualified by the laboratory with a "B" (indicative of method blank contamination) or an "A" (indicative of a common adol laboratory contaminant) were rejected "R".

<u>Fraction</u>	TIC	Samples
VOA	Column Bleed Hexane	CTS-05 CTS-03, -03RE, -04, -05, -06RE, -07, -07RE, -08, -08RE, -09, -09RE, -10, -10RE & -11RE

Note: TIC compounds associated with a "best match" spectra and CAS number were qualified as presumptive evidence of a compound at an estimated value "JN" by the data reviewer. TICs not associated with a CAS number were qualified as estimated "J" by the data reviewer in the sample data.

3. MASS SPECTROMETER TUNING:

Tuning and performance criteria are established to ensure adequate mass resolution, proper identification of compounds, and to some degree, sufficient instrument sensitivity. These criteria are not sample specific. Instrument performance is determined using standard materials. Therefore, these criteria should be met in all circumstances. The tuning standard for volatile organics is bromofluorobenzene (BFB) and for semi-volatiles is decafluorotriphenyl-phosphine (DFTPP).

If the mass calibration is in error or missing, all associated data will be classified as unusable "R". The following samples shown were qualified with "R" because of tuning:

VOA/BNA: Data met the QC criteria

4. CALIBRATION:

Satisfactory instrument calibration is established to ensure that the instrument is capable of producing acceptable quantitative data. An initial calibration demonstrates that the instrument is capable of giving acceptable performance at the beginning of an experimental sequence. The continuing calibration verifies that the instrument is giving satisfactory daily performance.

A) Response Factor:

The response factor measures the instrument's response to specific chemical compounds. The response factor for the VOA/BNA Target Compound List (TCL) must be ≥ 0.05 in both the initial and continuing calibrations. A value ≤ 0.05 indicates a serious detection and quantitation problem (poor sensitivity). If the mean RRF of the initial calibration or the continuing calibration has a response factor < 0.05 for any analyte, those analytes detected in environmental samples will be qualified as estimated "J". All non-detects for those compounds will be rejected "R". The following analytes in the samples shown were qualified because of response factor:

Initial Calibration

<u>VOA</u> - The following compounds were either qualified as estimated "J" (positive values only) or rejected "R" (non-detected "U" values only) in the associated samples because the Initial Calibration Mean RRF value is < 0.05:

Compound

Oualifier

Associated Sample(s)

Data met the QC criteria

BNA - The following compounds were either qualified as estimated "J" (positive values only) or rejected "R" (non-detected "U" values only) in the associated samples because the Initial Calibration Mean RRF value is < 0.05:

Compound

Oualifier

Associated Sample(s)

Data met the QC criteria

Continuing Calibration

 $\underline{\text{VOA}}$ - The following compounds were either qualified as estimated "J" (positive values only) or rejected "R" (non-detected "U" values only) in the associated samples because the Continuing Calibration RRF₅₀ is < 0.05:

Compound

Qualifier

Associated Sample(s)

Data met the QC criteria

BNA - The following compounds were either qualified as estimated "J" (positive values only) or rejected "R" (non-detected "U" values only) in the associated samples because the Continuing Calibration RRF_{∞} is < 0.05:

Compound

<u>Oualifier</u>

Associated Sample(s)

Data met the QC criteria

5. CALIBRATION:

B) PERCENT RELATIVE STANDARD DEVIATION (%RSD) AND PERCENT DIFFERENCE (%D):

Percent RSD is calculated from the initial calibration and is used to indicate the stability of the specific compound response factor over increasing concentration. Percent D compares the response factor of the continuing calibration check to the mean response factor (RRF) from the initial calibration. Percent D is a measure of the instrument's daily performance. Percent RSD must be < 30% and %D must be < 25%. A value outside of these QC limits indicates potential detection and quantitation errors. For these reasons, all positive results are flagged as estimated, "J"; and non-detects are flagged "UJ". If %RSD and/or %D grossly exceed QC criteria, non-detect data may be qualified "R".

For the PESTICIDE/PCB fraction, if %RSD exceeds 20% for all analytes except for the 2 surrogates (which must not exceed 30% RSD), qualify all associated positive results "J" and non-detects "UJ".

The following analytes in the samples shown were qualified for %RSD and %D:

Initial Calibration

<u>VOA</u> - Positive values of the following compounds were qualified as estimated "J" in the associated samples because the Initial Calibration %RSD is between 30-90% when the mean RRF is > 0.05:

Compound

Associated Sample(s)

Data met the QC criteria

BNA - Positive values of the following compounds were qualified as estimated "J" in the associated samples because the Initial Calibration %RSD is between 30-90% when the mean RRF is > 0.05:

Compound

Associated Sample(s)

Data met the QC criteria

Pest/PCBs - The following compounds were qualified as estimated "J" or rejected "R" in the associated samples because the linearity criteria or the percent relative standard deviation (%RSD) of the Initial Calibration is > 20% for either one or both GC columns:

Compound	Percent Recovery	Qualifier	Associated Sample(s)
Gamma BHC	22.7 %	"J"	CTS-01, -02, -03, -04, -05, -06, -07, -08, -9, -10, -11, CTW-01, -02, -03, -04, -05, -06, -07, -08,
4'4-DDT	27.5 %	"J"	-09 & -10 -01, -02, -03, -04, -05, -06, -07, -08, -9, -10, -11, CTW-01, -02, -03, -04, -05, -06, -07, -08, -09 & -10

Continuing Calibration

 \underline{VOA} - The following compounds were qualified as estimated "J" because the Continuing Calibration %D is between 25-90% when the RRF_{so} is > 0.05:

Compound	Associated Sample(s)
Acetone	CTS-01, -03, -06, -04, -06RE, -03RE, -08RE, -09RE, 10
	& -11RE
2-Butanone	CTS-04, -06RE, -03RE, -08RE, -09RE, 10 & -11RE
4-Methyl-2-Pentanone	CTS-04, -06RE, -03RE, -08RE, -09RE, 10 & -11RE
2-Hexanone	CTS-04, -06RE, -03RE, -08RE, -09RE, 10 & -11RE

<u>BNA</u> - The following compounds were qualified as estimated "J" because the Continuing Calibration %D is between 25-90% when the RRF₅₀ is > 0.05:

Compound	Associated Sample(s)
Hexachlorocyclopentadiene	CTS-03DL, -01DL, -08DL, -09DL, -10DL, -11DL & CTW-03
3-Nitroaniline	CTS-03DL, -01DL, -08DL, -09DL, -10DL & -11DL
Diethylphthalate	CTS-01, -02, -03, -04, -05, -06, -07, -08, -09, -10, -11, -4DL, -05DL, -06DL & -07DL

Compound	Associated Sample(s)
4-Chlorophenyl-phenylether	CTS-03DL, -01DL, -08DL, -09DL, -10DL, -11DL, CTS-01, -02, -03, -04, -05, -06, -07, -08, -09, -10, -11, 4DL, -05DL, -06DL & -07DL
4,6-Dinitro-2-methylphenol	CTS-01, -02, -03, -04, -05, -06, -07, -08, -09, -10, -11, -04DL, -05DL, -06DL & -07DL
Benzo (g.h.i)pervlene	CTW-06, -02RE, -08RE, -05RE, -07DL, -04DL, -01DL

09DL and -03

The following data were not qualified as estimated "J" in the associated samples because the percent difference (%D) of the Continuing Calibration is just outside specified QC Limits:

Fraction	Compound	OC Limits	Associated Sample(s)
BNA	4,6-Dinitro-2-Methylphenol(25.3%)	< 25.0%	CTS-03DL, -01DL, - 08DL, -09DL, -10DL &
			-11DL

Pest/PCBs - The Relative Percent Difference (%RSD) for PEM compound amounts in the continuing calibration verification analyses and/or the RPD amounts in the Individual Standard Mixes of the continuing calibration verification analyses are ≥ 25% for either one or both GC columns. The following compounds were either qualified as estimated "J" or rejected "R" due to exceeding Continuing Calibration QC criteria:

Compound RPD Qualifier Associated Sample(s)

Data met QC criteria

Associated samples are those samples located approximately five (5) samples on either side of the failing PEM sample and/or between the last in-control standard and the next passing standard.

Pest/PCBs - The following compounds were qualified as estimated "J" in the associated samples because the Continuing Calibration %D is between 20-90% for these compounds on the primary GC column:

Compound

Associated Samples

Data met QC criteria

6. SURROGATES/SYSTEM MONITORING COMPOUNDS (SMC):

All samples are spiked with surrogate/SMC compounds prior to sample preparation to evaluate overall laboratory performance and efficiency of the analytical technique. If the measured surrogate/SMC concentrations were outside contract specifications, qualifications were applied to the samples and analytes as shown below. The following analytes for the samples shown were qualified because of surrogate/SMC recovery:

<u>VOA</u> - The following compounds were either qualified as estimated "J" or rejected "R" due to surrogate recovery outside specified QC limits:

Surrogate

Recovery

Oualifier

Compounds

Sample(s)

Data met the QC criteria.

BNA - The following compounds were either qualified as estimated "J" or rejected "R" due to surrogate recovery outside specified QC limits:

Surrogate

Recovery

Oualifier

Compounds

Sample(s)

Data met the QC criteria.

Note: Data were qualified because either two (2) base-neutral and/or two (2) acid compounds have recoveries outside specified QC limits and above 10%, or either one (1) base-neutral and/or one (1) acid compound has a percent recovery below 10%.

Pest/PCBs - The following compounds were either qualified as estimated "J" or rejected "R" due to Tetrachloro-m-xylene (TCX) and Decachlorobiphenyl (DCB) surrogate recoveries are both outside specified advisory QC limits (60-150%):

Surrogate	Recovery	Qualifier	Compounds	Sample(s)
Decachlorobiphenyl (DCB)	< 10%	"R" "J ¹ "	25 03	CTS-04 CTS-04

Positive values only were qualified as estimated "J" in the associated samples.

Surrogate recoveries outside QC criteria (< 10%) may be attributable to the required dilution of the extract during analysis and/or due to matrix interference.

Note: Data were qualified because recoveries for both surrogates are outside specified QC limits and above 10%, or either surrogate has a percent recovery below 10%.

7. INTERNAL STANDARDS PERFORMANCE:

Internal standard (IS) performance criteria ensure that the GC/MS sensitivity and response are stable during every experimental run. The internal standard area count must not vary by more than a factor of 2 (-50% to 100%) from the associated continuing calibration standard. The retention time of the internal standard must not vary more than \pm 30 seconds from the associated continuing calibration standard. If the area count is outside the -50% to 100% range of the associated standard, all of the positive results for compounds quantitated using that IS are qualified as estimated "J", and all non-detects as "UJ"only if the IS area is <50% Non-detects are qualified as "R" if there is a severe loss of sensitivity (<25% of associated IS area counts).

If an internal standard retention time varies by more than 30 seconds, the reviewer will use professional judgement to determine either partial or total rejection of the data for that sample fraction. The following analytes in the samples shown were qualified because of internal standard performance:

<u>VOA</u> - The following compounds were either qualified as estimated "J" or rejected "R" in the associated samples due to exceeding Internal Standard (IS) QC criteria (within -50% to + 100% of the Continuing Calibration 12-hour standard):

			•
Internal Standard	Percent IS Area Count of the 12-Hour Standard	Total Analytes Qualified/Sample	Associated Sample(s)
Bromochloromethane	between -25% to -50%	"J"/13	CTS-03, -03RE, -08RE, -10, -11RE, -08, -10RE, -11,*-07RE & -07
1,4 - Difluorobenzene	between -25% to -50%	"J"/11	CTS-03, -03RE, -08RE, -10, -11RE, -08, -10RE, -11, -07RE & -07
Chlorobenzene-d5	between -25% to -50%	"J"/9	CTS-06, -06RE, -09RE, -10, -09, -10RE, -11, -02
	< -25%	"R"/9	CTS-03, -03RE, 08RE, - 11RE, -08, -07RE & -07

^{*} The sample CTS-07RE was not rejected as "R" because the IS area count of the 12-Hour standard is just below the <-25%:

Positive values only were qualified as estimated "J" in the indicated samples.

7. INTERNAL STANDARDS PERFORMANCE (continued):

BNA - The following compounds were either qualified as estimated "J" or rejected "R" in the associated samples due to exceeding Internal Standard (IS) QC criteria (within -50% to + 100% of the Continuing Calibration 12-hour standard):

Internal Standard	Percent IS Area Count of the 12-Hour Standard	Qualifier/Total Analytes Qualified	Associated Sample(s)
Chrysene-d12	between -25% to -50%	"J"/6	CTS-01, -04 & 11 CTW-07, -04, -02, -08, - 05 & -09
	< -25%	"R"/7	CTW-01
Perylene-d12	between -25% to -50%	"J"	CTS-05, -06, -08, -09, - 10, -02 & -04DL, -07DL CTW-10, 02RE, -08RE
	< -25%	"R"/7	& -05RE CTS-01, -03, -04, -07 &
			CTW-07, -04, -01, -02, - 08, -05, -09, -07DL, - 04DL, -01DL & -09DL

Note: The laboratory indicated in the case narrative that samples exhibited internal standard areas outside QC criteria due to a matrix effect confirmed by repeat analysis.

INTERNAL STANDARD AREA OUTLIERS

Sample	Internal Standard VOA Fraction	Area	Lower Limit	Upper Limit	< -25% of the 12-hour Standard
CTS-03	всм	15236	18213	72852	9106.5
CTS-03	DFB	67158	85114	340456	42557
CTS-03	CBZ	32936	68723	274892	34361.5
CTS-06	CBZ	57533	68723	274892	34361.5
CBZ-06RE	CBZ	67271	72111	288442	36055.5
CTS-03RE	ВСМ	11366	19485	77940	9742.5
CTS-03RE	DFB	52740	90139	360556	45069.5
CTS-03RE	CBZ	22779	72111	288442	36055.5
CTS-08RE	ВСМ	12499	19485	77940	9742.5
CTS-08RE	DFB	49892	90139	360556	45069.5
CTS-08RE	CBZ	20655	72111	288442	36055.5
CTS-09RE	CBZ	56670	72111	288442	36055.5
CTS-10	ВСМ	17976	19485	77940	9742.5
CTS-10	DFB	77973	90139	360556	45069.5
CTS-10	CBZ	48699	72111	288442	36055.5
CTS-11RE	BCM	13829	19485	77940	9742.5
	DFB	59374	90139	360556	45069.5
CTS-11RE	CBZ	35636	72111	288442	36055.5
CTS-11RE	BCM	12153	17788	71152	8894
CTS-08	DFB	57111	81086	324344	40543
CTS-08	CBZ	32897	67809	271234	33904.5
CTS-08	CBZ	56066	67809	271234	33904.5
CTS-10PF	ВСМ	17050	17788	71152	8894
CTS-10RE	DFB	78294	81086	324344	40543
CTS-10RE	CBZ	45482	67809	271234	33904.5
CTS-10RE		17700	17788	71152	8894
CTS-11	BCM	78945	81086	324344	40543
CTS-11	DFB		67809	271234	33904.5
CTS-11	CBZ	51372	67809	271234	33904.5
CTS-02	CBZ	60195	07809	211237	

Sample	Internal Standard VOA Fraction	Area	Lower Limit	Upper Limit	< -25% of the 12-hour Standard
CTS-07RE	ВСМ	8893	17788	71152	8894
CTS-07RE	DFB	42749	81086	324344	40543
CTS-07RE	CBZ	13597	67809	271234	33904.5
CTS-07	BCM	12878	17788	71152	8894
	DFB	55406	81086	324344	40543
CTS-07	CBZ	22451	67809	271234	33904.5

INTERNAL STANDARD AREA OUTLIERS

Sample	Internal Standard BNA Fraction	Area	Lower Limit	Upper Limit	< -25% of the 12-hour Standard
CTS-01	CRY	29886	38566	154264	19283
CTS-01	PRY	9624	34691	138762	17345.5
CTS-03	PRY	10894	34691	138762	17345.5
CTS-04	CRY	31024	38566	154264	19283
CTS-04	PRY	13884	34691	138762	17345.5
CTS-05	PRY	22541	34691	138762	17345.5
CTS-06	PRY	23137	34691	138762	17345.5
CTS-07	PRY	15933	34691	138762	17345.5
CTS-08	PRY	25672	34691	138762	17345.5
CTS-09	PRY	24925	34691	138762	17345.5
CTS-10	PRY	18176	34691	138762	17345.5
CTS-11	CRY	27741	38566	154264	19283
CTS-11	PRY	11466	34691	138762	17345.5
CTS-02	PRY	32355	34691	138762	17345.5
CTS-04DL	PRY	33013	34691	138762	17345.5
CTS-07DL	PRY	30710	34691	138762	17345.5
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INTERNAL STANDARD AREA OUTLIERS

Sample	Internal Standard BNA Fraction	Area	Lower Limit	Upper Limit	< -25% of the 12-hour Standard
CTW-10	PRY	19789	26387	105548	13194
CTW-07	CRY	22065	34616	138462	17308
CTW-07	PRY	13172	26387	105548	13194
CTW-04	CRY	20261	34616	138462	17308
CTW-04	PRY	13099	26387	105548	13194
CTW-01	CRY	16485	34616	138462	17308
CTW-01	PRY	10925	26387	105548	13194
CTW-02	CRY	18812	34616	138462	17308
CTW-02	PRY	10719	26387	105548	13194
CTW-08	CRY	17549	34616	138462	17308
CTW-08	PRY	10516	26387	105548	13194
CTW-05	CRY	18439	34616	138462	17308
CTW-05	PRY	11877	26387	105548	13194
CTW-09	CRY	17351	34616	138462	17308
CTW-09	PRY	11219	26387	105548	13194
CTW-02RE	PRY	37963	51133 .	204532	25566
CTW-02RE	PRY	35375	51133	204532	25566
CTW-05RE	PRY	44533	51133	204532	25566
CTW-07DL	PRY	20511	51133	204532	25566
CTW-04DL	PRY	21285	51133	204532	25566
CTW-01DL	PRY	22823	51133	204532	25566
CTW-09DL	PRY	21890	51133	204532	25566
C1 11-07DD					

8. COMPOUND IDENTIFICATION:

A) VOLATILE AND SEMI-VOLATILE FRACTIONS:

TCL compounds are identified on the GC/MS by using the analyte's relative retention time (RRT) and by comparison to the ion spectra obtained from known standards. For the results to be a positive hit, the sample peak must be within \pm 0.06 RRT units of the standard compound, and have an ion spectra which has a ratio of the primary and secondary m/e intensities within 20% of that in the standard compound. For the Tentatively Identified Compounds (TICs) the ion spectra must match accurately. In the cases where there is not an adequate ion spectrum match, the laboratory may have provided false positive identifications. The following analytes in the samples shown were qualified for compound identification:

The following compounds were qualified as estimated "J" in the indicated samples because they could not be chromatographically resolved:

Fraction

Compounds

Samples

Data met the QC criteria.

B) PESTICIDE FRACTION:

The retention time of the reported compounds must fall within the calculated retention time windows for the two chromatographic columns and a GC/MS confirmation is required if the concentration exceeds 10 ng/ml in the final sample extract. The percent difference (%D) of the positive results obtained on the two GC columns would be $\leq 25\%$. The following analytes in the samples shown were qualified because of compound identification:

<u>Pest/PCBs</u> - The following detected compounds were qualified due to a percent difference (%D) between the primary and confirmation columns > 25%:

Compound	<u>%D</u>	Qualifier	Sample(s)
beta-BHC	between 25-50% between 50-90%	"J"	CTS-03, -06, -07 CTS-04
delta-BHC	between 50-90%	"JN"	CTS-06
4,4'-DDD	between 50-90%	"JN"	CTS-02, -04, -06
4,4'-DDT	between 50-90%	"JN"	CTS-06
Aroclor-1254	between 25-50% between 50-90%	"J" "JN"	CTS-02, -04 ¹ , -10 & CTW-04 CTS-08

This sample was previously qualified as estimated "J" due to surrogate recovery criteria.

<u>Pest/PCBs</u> - Due to professional judgement, the lower of two positive values generated by the laboratory from the primary and confirmation column analyses was used to report final results for the following pesticide compounds:

Compound

Primary Column Value Confirmation Column Value

No qualification was required based on this criteria.

Note: During the initial calibration sequence, absolute retention times are determined for all single response pesticides, the surrogates, and at least three major peaks of each multi-component analyte. Windows are centered around the mean absolute retention time for the analyte established during the initial calibration. Analytes are identified when peaks are observed in the retention time window for the compound on both GC columns. The quant reports listed many potential pesticide compounds for consideration. Comparison of the sample retention times to the retention time windows established during the initial calibration revealed that no additional pesticide compounds were detected in the associated samples. In addition, no shifts for surrogate compound retention times were noted to occur that might require consideration of compounds outside respective retention time windows.

9. MATRIX SPIKE/SPIKE DUPLICATE, MS/MSD:

The MS/MSD data are generated to determine the long-term precision and accuracy of the analytical method in various matrices. The MS/MSD may be used in conjunction with other QC criteria for some additional qualification of the data. The following analytes, for the samples shown, were qualified because of MS/MSD:

The laboratory indicated in the case narrative that sample CTS-02 for soil samples, CTW-10 for wipe samples and CTS-05 for med-conc. VOA samples were used as the original to prepare the duplicate matrix spikes.

<u>VOA</u> - The following sample data were either qualified as estimated "J" or rejected "R" due to exceeding duplicate spike recovery QC criteria:

Original Sample	Spike Recovery	Qualifier		Compound(s)
CTS-05(med-conc.)	< 10%	"J 1"	•'	Toluene

Previosly qualified due to holding time criteria.

BNA - The following sample data were either qualified as estimated "J" or rejected "R" due to exceeding duplicate spike recovery QC criteria:

Original Sample	Spike Recovery	Qualifier	Compound(s)
CTS-02	< 10%	"J¹"	Pyrene

Positive values only were qualified as estimated "J" in the indicated samples.

<u>Pest/PCBs</u> - The following sample data were either qualified as estimated "J" or rejected "R" due to exceeding duplicate spike recovery QC criteria:

Original Sample Spike Recovery Qualifier Compound(s)

No qualification was required.

10. OTHER QC DATA OUT OF SPECIFICATION:

Full TCL - The following compounds were qualified as estimated "J" in the associated aqueous and/or soil/sediment field duplicate samples because the Relative Percent Difference (RPD) between the sample and field duplicate sample is > 50% for aqueous samples, or > 100% for soil/sediment samples:

Compound

Matrix

% RPD

Associated Field Duplicate Samples

Data met the QC criteria.

The following soil/sediment/solid sample data (other than TCLP data) were either qualified as estimated "J" (% solids between 10-50%) or rejected "R" (% solids < 10%) because the sample contains more than 50% water:

Fraction

Percent Solids

Oualifier

Compounds Sample(s)

Data met the QC criteria.

<u>VOA</u> - The following data were either qualified as estimated "J" or rejected "R" due to air bubbles in the VOA vial(s):

Sample(s)

Oualifier

Compounds

No qualification was required.

10. OTHER QC DATA OUT OF SPECIFICATION (continued):

The following compounds were qualified as estimated "J" in the indicated samples because the on-column amount of these compounds exceeded the instrument's analytical range as defined by the highest concentration level of the Initial Calibration Sequence:

Fraction	Sample(s)	Compound(s)
BNA	bis(2-Ethylhexyl)phthalate	CTS-01 ¹ , -03, -05, -07, -08, -09, -10 & 11 ¹ CTW-07 ¹ , -09 ¹ & -01 ¹
	Benzo(a)anthracene Chrysene Benzo(b)Fluranthene Benzo(a)Pyrene Pyrene Butylbenzylphthalate	CTS-04 ¹ CTS-04 ¹ CTS-04 ¹ CTS-04 ¹ CTS-04 ¹ CTS-04 ¹ & CTS-06 CTS-05, -08 & -10 CTW-07 ¹ , -09 ¹ , -01 ¹ & -04 ¹
	Di-n-octylphthalate	CTS-10 ¹
PCBs	Aroclor-1254	CTS-01, -03, -05, -07 & 11

These samples were previously qualified as estimated "J" due to other QC criteria.

11. SYSTEM PERFORMANCE AND OVERALL ASSESSMENT:

Due to professional judgement, the following compounds were not transferred from the indicated dilution sample analyses to the undiluted sample analyses because the reported values of these compounds are either diluted out in the associated dilution sample analyses or are qualified as non-detected "U" due to blank contamination QC criteria:

_		
Fra	ction	

Compound

Dilution Sample(s)

Dilution Factor

No qualification was required based on this criteria.

Due to professional judgement, the following positive data were rejected "R" due to possible carryover from a previous sample analysis that contained the compound(s) at high concentration(s):

Fraction

Sample Compound

Sample Compound

Previous Sample Compound Concentration

Concentration

No qualification was required based on this criteria.

12. CONTRACT PROBLEMS NON-COMPLIANCE:

VOA: VOA analysis was not performed on wipe samples due to analytical problems with hexane preservatives.

BNA:

- 1) GPC Clean-up was not performed for wipe samples.
- 2) Labs failed to submit undiluted results for the sample CTW-03.
- 3) Internal Standard Perylene-d12 is below the lower limit in sample CTW-10, laboratory failed to reanalyze the sample.

PCBs: A multi component analyte Aroclor - 1254 was detected in most of the samples, but a matching multi component standard was not analyzed within 72 hours of the injection of the sample and within a valid 12 hour sequence. No qualification was required based on this criteria.

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In Association with Resource Applications, Inc., R.E. Serviera Associates, PRC Environmental Management, C.C. Johnson & Malhotra, P.C., and GRB Environmental Services, Inc.

START DATA SIGN-OFF SHEET

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